

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN032719\
 Data File : BN004992.D
 Acq On : 27 Mar 2019 10:49
 Operator : JU/SJ
 Sample : K1825-03
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-BN-WP

Quant Time: Mar 29 04:36:05 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-BN032619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 27 05:27:09 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	9782	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	39340	20.00	ng	0.00
39) Acenaphthene-d10	14.32	164	21973	20.00	ng	0.00
64) Phenanthrene-d10	17.07	188	48256	20.00	ng	0.00
76) Chrysene-d12	21.27	240	46964	20.00	ng	0.00
87) Perylene-d12	23.51	264	45735	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	79339	140.23	ng	0.00
7) Phenol-d6	6.86	99	105704	140.86	ng	0.00
23) Nitrobenzene-d5	8.85	82	62941	96.78	ng	0.00
42) 2,4,6-Tribromophenol	15.82	330	42296	119.35	ng	0.00
45) 2-Fluorobiphenyl	12.94	172	163190	98.42	ng	0.00
79) Terphenyl-d14	19.70	244	244424	100.20	ng	0.00

Target Compounds				Qvalue		
4) n-Nitrosodimethylamine	3.52	42	12838	68.345	ng	96
12) 1,3-Dichlorobenzene	7.56	146	60667	77.525	ng	99
13) 1,4-Dichlorobenzene	7.71	146	90742	114.558	ng	100
16) 2,2'-oxybis(1-Chloropropan	8.20	45	91170	121.522	ng	100
18) Hexachloroethane	8.74	117	33775	120.299	ng	99
19) n-Nitroso-di-n-propylamine	8.48	70	38513	78.286	ng	98
24) Nitrobenzene	8.89	77	57544	87.431	ng	99
28) bis(2-Chloroethoxy)methane	9.88	93	24177	30.800	ng	99
30) 1,2,4-Trichlorobenzene	10.32	180	71026	105.441	ng	99
31) Naphthalene	10.51	128	167477	81.596	ng	100
49) Acenaphthylene	14.05	152	191281	79.690	ng	100
50) Dimethylphthalate	13.79	163	184642	102.281	ng	99
51) 2,6-Dinitrotoluene	13.92	165	48152	117.221	ng	97
52) Acenaphthene	14.39	154	20215	14.971	ng	100
57) 2,4-Dinitrotoluene	14.70	165	40116	71.860	ng	97
60) Diethylphthalate	15.15	149	196986	109.351	ng	99
61) 4-Chlorophenyl-phenylether	15.37	204	62053	69.303	ng	98
66) n-Nitrosodiphenylamine	15.59	169	74301	48.881	ng	99
67) 4-Bromophenyl-phenylether	16.26	248	35032	58.175	ng	98
74) Di-n-butylphthalate	18.03	149	184109	61.523	ng	100
75) Fluoranthene	19.15	202	442539	143.064	ng	99
80) Butylbenzylphthalate	20.40	149	77668	60.954	ng	99
81) Benzo(a)anthracene	21.26	228	400520	136.912	ng	99
83) Chrysene	21.32	228	368438	131.857	ng	99
84) Bis(2-ethylhexyl)phthalate	21.17	149	268922	152.537	ng	98
85) Di-n-octyl phthalate	22.05	149	125898	44.249	ng	99
90) Benzo(a)pyrene	23.42	252	146154	56.734	ng	99
92) Benzo(g,h,i)perylene	26.49	276	340722	133.215	ng	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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