

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP111119\
 Data File : BP001032.D
 Acq On : 11 Nov 2019 23:09
 Operator : JU
 Sample : K5676-11
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SU-03-110819

Manual Integrations
 APPROVED

mohammad
 11/12/2019 5:04:51 PM

Quant Time: Nov 12 06:50:02 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP110619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 06 17:07:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	235913	20.00	ng	0.00
21) Naphthalene-d8	10.40	136	852207	20.00	ng	0.00
39) Acenaphthene-d10	13.26	164	476890	20.00	ng	0.00
64) Phenanthrene-d10	15.66	188	887575	20.00	ng	0.00
76) Chrysene-d12	19.30	240	693233	20.00	ng	0.00
87) Perylene-d12	21.09	264	665963	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.31	112	1221058	94.21	ng	0.01
7) Phenol-d6	7.82	99	1604667	97.91	ng	0.00
23) Nitrobenzene-d5	9.25	82	985945	64.00	ng	0.00
42) 2,4,6-Tribromophenol	14.56	330	574247	100.87	ng	0.00
45) 2-Fluorobiphenyl	12.18	172	2048880	64.71	ng	0.00
79) Terphenyl-d14	17.94	244	2302716	66.07	ng	0.00

Target Compounds

						Qvalue
10) Phenol	7.83	94	104294	5.818	ng	98
49) Acenaphthylene	13.03	152	169520	3.885	ng	100
50) Dimethylphthalate	12.85	163	182008	5.161	ng	100
52) Acenaphthene	13.31	154	69904	2.678	ng	94
58) Fluorene	14.16	166	111280	3.430	ng	99
71) Phenanthrene	15.69	178	1357051	29.873	ng	100
72) Anthracene	15.77	178	393928	8.889	ng	99
73) Carbazole	16.01	167	101539	2.499	ng	98
75) Fluoranthene	17.40	202	2198479	42.817	ng	99
78) Pyrene	17.71	202	2022583	41.285	ng	98
81) Benzo(a)anthracene	19.29	228	1046486	22.414	ng	94
83) Chrysene	19.33	228	1017005	24.811	ng	97
86) Indeno(1,2,3-cd)pyrene	22.76	276	409904	7.309	ng	# 92
88) Benzo(b)fluoranthene	20.59	252	1145464m	28.855	ng	
89) Benzo(k)fluoranthene	20.62	252	308448m	8.254	ng	
90) Benzo(a)pyrene	21.02	252	740781	20.294	ng	95
91) Dibenzo(a,h)anthracene	22.78	278	107485	2.885	ng	# 86
92) Benzo(g,h,i)perylene	23.25	276	356326	9.499	ng	96

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

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