

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP010320\
 Data File : BP001513.D
 Acq On : 03 Jan 2020 18:24
 Operator : JU
 Sample : K6471-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MC-123019-0-2-COMP-B5

Quant Time: Jan 04 03:16:03 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 03 15:57:46 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	77863	20.00	ng	0.00
21) Naphthalene-d8	10.48	136	285810	20.00	ng	0.00
39) Acenaphthene-d10	14.35	164	171837	20.00	ng	0.00
64) Phenanthrene-d10	17.10	188	370514	20.00	ng	0.00
76) Chrysene-d12	21.20	240	323590	20.00	ng	0.00
87) Perylene-d12	23.46	264	347877	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.35	112	449189	96.92	ng	0.04
7) Phenol-d6	6.93	99	601878	96.42	ng	0.05
23) Nitrobenzene-d5	8.86	82	349905	60.47	ng	0.01
42) 2,4,6-Tribromophenol	15.85	330	170613	87.73	ng	0.00
45) 2-Fluorobiphenyl	12.97	172	679300	58.13	ng	0.00
79) Terphenyl-d14	19.69	244	877713	54.06	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	6.95	94	18717	2.850	ng	94
50) Dimethylphthalate	13.83	163	38002	2.829	ng	99
71) Phenanthrene	17.15	178	54836	2.729	ng	98
74) Di-n-butylphthalate	18.10	149	48869	2.089	ng	97
75) Fluoranthene	19.12	202	134611	5.533	ng	98
78) Pyrene	19.47	202	130507	5.558	ng	99
81) Benzo(a)anthracene	21.18	228	64933	2.860	ng	98
83) Chrysene	21.23	228	68745	3.158	ng	99
84) Bis(2-ethylhexyl)phthalate	21.16	149	42130	2.846	ng	# 97
88) Benzo(b)fluoranthene	22.78	252	86835	4.031	ng	# 83
90) Benzo(a)pyrene	23.36	252	63853	3.167	ng	# 85
92) Benzo(a,h,i)perylene	26.43	276	46326	2.318	ng	# 89

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

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