

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
 Data File : BP001374.D
 Acq On : 23 Dec 2019 18:38
 Operator : JU
 Sample : K6401-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B38

Quant Time: Dec 24 05:51:14 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 20 03:20:39 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.30	152	52442	20.00	ng	0.00
21) Naphthalene-d8	10.34	136	210473	20.00	ng	0.00
39) Acenaphthene-d10	13.20	164	112434	20.00	ng	0.00
64) Phenanthrene-d10	15.60	188	217410	20.00	ng	0.00
76) Chrysene-d12	19.25	240	148321	20.00	ng	0.00
87) Perylene-d12	21.03	264	150708	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	6.22	112	299702	92.36	ng	0.02
7) Phenol-d6	7.76	99	418448	98.83	ng	0.01
23) Nitrobenzene-d5	9.19	82	291047	53.10	ng	0.00
42) 2,4,6-Tribromophenol	14.50	330	133053	94.65	ng	0.00
45) 2-Fluorobiphenyl	12.12	172	518093	58.27	ng	0.00
79) Terphenyl-d14	17.89	244	588813	67.99	ng	0.00
Target Compounds						
10) Phenol	7.78	94	14088	2.912	ng	84
17) 2-Methylphenol	8.73	107	90970	31.558	ng	97
20) 3+4-Methylphenols	8.99	107	463055	121.052	ng	92
27) 2,4-Dimethylphenol	9.83	122	150647	53.248	ng	98
50) Dimethylphthalate	12.79	163	82994	8.998	ng	99
84) Bis(2-ethylhexyl)phthalate	19.29	149	34710	4.406	ng	# 96

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

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