

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
 Data File : BP001385.D
 Acq On : 24 Dec 2019 14:35
 Operator : JU
 Sample : K6360-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 1FT-TP

Quant Time: Dec 24 16:20:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 16:04:49 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.28	152	50826	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	187621	20.00	ng	0.00
39) Acenaphthene-d10	13.17	164	113492	20.00	ng	0.00
64) Phenanthrene-d10	15.57	188	220071	20.00	ng	0.00
76) Chrysene-d12	19.22	240	149222	20.00	ng	0.00
87) Perylene-d12	20.99	264	140366	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.21	112	362299	115.21	ng	0.01
7) Phenol-d6	7.75	99	506289	123.38	ng	0.00
23) Nitrobenzene-d5	9.17	82	348604	71.35	ng	0.00
42) 2,4,6-Tribromophenol	14.47	330	168194	118.54	ng	0.00
45) 2-Fluorobiphenyl	12.10	172	674142	75.11	ng	0.00
79) Terphenyl-d14	17.87	244	786505	90.27	ng	0.00
Target Compounds						
10) Phenol	7.77	94	11684	2.492	ng	Qvalue # 81
50) Dimethylphthalate	12.76	163	33396	3.587	ng	98
71) Phenanthrene	15.61	178	26804	2.135	ng	96
75) Fluoranthene	17.32	202	67713	4.823	ng	98
78) Pyrene	17.63	202	51149	4.415	ng	97
83) Chrysene	19.25	228	22692	2.162	ng	96
88) Benzo(b)fluoranthene	20.50	252	25780	2.797	ng	# 96

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

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