

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP111119\
 Data File : BP001031.D
 Acq On : 11 Nov 2019 22:39
 Operator : JU
 Sample : K5778-07
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-35-S

Quant Time: Nov 12 06:45:58 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	237826	20.00	ng	0.00
21) Naphthalene-d8	10.40	136	857201	20.00	ng	0.00
39) Acenaphthene-d10	13.26	164	494936	20.00	ng	0.00
64) Phenanthrene-d10	15.66	188	977649	20.00	ng	0.00
76) Chrysene-d12	19.29	240	806239	20.00	ng	0.00
87) Perylene-d12	21.08	264	827565	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.31	112	1265598	96.86	ng	0.01
7) Phenol-d6	7.82	99	1652952	100.04	ng	0.00
23) Nitrobenzene-d5	9.25	82	1028481	66.37	ng	0.00
42) 2,4,6-Tribromophenol	14.56	330	620782	105.07	ng	0.00
45) 2-Fluorobiphenyl	12.18	172	2116541	64.41	ng	0.00
79) Terphenyl-d14	17.94	244	2588336	63.86	ng	0.00
Target Compounds						
10) Phenol	7.83	94	103473	5.725	ng	95
50) Dimethylphthalate	12.85	163	145035	3.963	ng	100
78) Pyrene	17.70	202	138317	2.428	ng	99
88) Benzo(b)fluoranthene	20.58	252	122913	2.492	ng	95
90) Benzo(a)pyrene	21.00	252	95264	2.100	ng	97

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

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