

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP111519\
 Data File : BP001057.D
 Acq On : 15 Nov 2019 23:56
 Operator : JU
 Sample : K5832-17
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 5-(6-8)

Manual Integrations
 APPROVED

mohammad
 11/18/2019 3:06:40 PM

Quant Time: Nov 16 04:56:56 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	8.37	152	199744	20.00	ng	0.00
21) Naphthalene-d8	10.39	136	715696	20.00	ng	-0.01
39) Acenaphthene-d10	13.25	164	414097	20.00	ng	-0.02
64) Phenanthrene-d10	15.65	188	806515	20.00	ng	-0.01
76) Chrysene-d12	19.28	240	663962	20.00	ng	-0.02
87) Perylene-d12	21.06	264	714103	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	6.31	112	1258636	114.69	ng	0.01
7) Phenol-d6	7.82	99	1726512	124.41	ng	0.00
23) Nitrobenzene-d5	9.25	82	1098883	84.93	ng	-0.01
42) 2,4,6-Tribromophenol	14.55	330	595397	120.44	ng	-0.01
45) 2-Fluorobiphenyl	12.17	172	2346431	85.34	ng	-0.01
79) Terphenyl-d14	17.93	244	2885734	86.45	ng	-0.01
Target Compounds						
10) Phenol	7.83	94	56016	3.690	ng	90
50) Dimethylphthalate	12.84	163	387785	12.664	ng	99
71) Phenanthrene	15.68	178	100236	2.428	ng	98
75) Fluoranthene	17.39	202	208765	4.474	ng	99
78) Pyrene	17.69	202	235637	5.022	ng	98
81) Benzo(a)anthracene	19.27	228	114446	2.559	ng	93
83) Chrysene	19.32	228	136422	3.475	ng	98
88) Benzo(b)fluoranthene	20.57	252	172086m	4.043	ng	
90) Benzo(a)pyrene	20.99	252	88065	2.250	ng	99

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

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