

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP111519\
 Data File : BP001058.D
 Acq On : 16 Nov 2019 00:26
 Operator : JU
 Sample : K5832-18
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 5-(8-10)

Manual Integrations
 APPROVED

mohammad
 11/18/2019 3:06:43 PM

Quant Time: Nov 16 05:09:36 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	163460	20.00	ng	0.00
21) Naphthalene-d8	10.39	136	608996	20.00	ng	-0.01
39) Acenaphthene-d10	13.25	164	375917	20.00	ng	-0.02
64) Phenanthrene-d10	15.65	188	775628	20.00	ng	-0.01
76) Chrysene-d12	19.28	240	664630	20.00	ng	-0.02
87) Perylene-d12	21.06	264	726914	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	6.30	112	526952	58.68	ng	0.00
7) Phenol-d6	7.81	99	922182	81.20	ng	0.00
23) Nitrobenzene-d5	9.24	82	609157	55.33	ng	-0.02
42) 2,4,6-Tribromophenol	14.55	330	220319	49.10	ng	-0.01
45) 2-Fluorobiphenyl	12.17	172	1290902	51.72	ng	-0.02
79) Terphenyl-d14	17.92	244	1759944	52.67	ng	-0.02
Target Compounds						
10) Phenol	7.83	94	30188	2.430	ng	95
50) Dimethylphthalate	12.83	163	232411	8.361	ng	99
71) Phenanthrene	15.68	178	384952	9.697	ng	99
72) Anthracene	15.76	178	78571	2.029	ng	97
75) Fluoranthene	17.39	202	381672	8.506	ng	98
78) Pyrene	17.69	202	518056	11.030	ng	98
81) Benzo(a)anthracene	19.27	228	217595	4.861	ng	94
83) Chrysene	19.32	228	236894	6.028	ng	98
88) Benzo(b)fluoranthene	20.57	252	246792m	5.696	ng	
90) Benzo(a)pyrene	20.99	252	163192	4.096	ng	98
92) Benzo(g,h,i)perylene	23.20	276	95695	2.337	ng	96

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

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