

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP100719\
 Data File : BP000549.D
 Acq On : 07 Oct 2019 17:43
 Operator : JU
 Sample : K5213-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 S-2(0-5)

Manual Integrations
 APPROVED

mohammad
 10/8/2019 12:23:57 PM

Quant Time: Oct 07 18:29:18 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 01 18:03:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	274467	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	987703	20.00	ng	-0.01
39) Acenaphthene-d10	9.80	164	543921	20.00	ng	-0.01
64) Phenanthrene-d10	11.30	188	989444	20.00	ng	0.00
76) Chrysene-d12	13.97	240	855460	20.00	ng	0.00
87) Perylene-d12	15.45	264	974341	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	1374432	80.49	ng	-0.01
7) Phenol-d6	6.40	99	2124767	103.27	ng	0.00
23) Nitrobenzene-d5	7.32	82	1163488	71.95	ng	-0.01
42) 2,4,6-Tribromophenol	10.60	330	362662	62.57	ng	-0.01
45) 2-Fluorobiphenyl	9.13	172	2194659	65.28	ng	0.00
79) Terphenyl-d14	12.91	244	2509842	59.40	ng	0.00
Target Compounds						
10) Phenol	6.41	94	86156	4.090	ng	98
50) Dimethylphthalate	9.52	163	339291	8.977	ng	100
71) Phenanthrene	11.33	178	189979	3.823	ng	100
75) Fluoranthene	12.53	202	355845	6.376	ng	99
78) Pyrene	12.76	202	362481	6.144	ng	99
81) Benzo(a)anthracene	13.96	228	200357	3.839	ng	95
83) Chrysene	14.00	228	167426	3.268	ng	97
88) Benzo(b)fluoranthene	15.02	252	232236m	4.201	ng	
90) Benzo(a)pyrene	15.39	252	169904	3.297	ng	95
92) Benzo(g,h,i)perylene	17.36	276	110260	2.171	ng	95

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

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