

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF013019\
 Data File : BF112324.D
 Acq On : 30 Jan 2019 16:23
 Operator : JU/SJ
 Sample : K1282-09
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3

Manual Integrations
 APPROVED

mohammad
 2/1/2019 12:06:16 AM

Quant Time: Jan 30 23:59:33 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 28 10:38:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	194871	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	734883	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	338585	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	587168	20.00	ng	0.00
76) Chrysene-d12	14.00	240	410169	20.00	ng	-0.01
87) Perylene-d12	15.47	264	414417	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	1001645	109.18	ng	0.00
7) Phenol-d6	6.49	99	1230773	96.54	ng	0.00
23) Nitrobenzene-d5	7.40	82	780522	61.20	ng	-0.01
42) 2,4,6-Tribromophenol	10.67	330	324728	82.40	ng	-0.01
45) 2-Fluorobiphenyl	9.19	172	1261369	52.69	ng	-0.01
79) Terphenyl-d14	12.94	244	973001	44.68	ng	-0.01
Target Compounds						
10) Phenol	6.50	94	91934	6.573	ng	80
50) Dimethylphthalate	9.59	163	155788	5.922	ng	97
71) Phenanthrene	11.38	178	198922	6.516	ng	97
75) Fluoranthene	12.57	202	275073	8.402	ng	97
78) Pyrene	12.80	202	236261	8.320	ng	98
81) Benzo(a)anthracene	13.99	228	118789	4.927	ng	97
83) Chrysene	14.03	228	120172	5.221	ng	95
86) Indeno(1,2,3-cd)pyrene	16.94	276	52860	2.898	ng	94
88) Benzo(b)fluoranthene	15.04	252	144305m	5.822	ng	
89) Benzo(k)fluoranthene	15.06	252	56391m	2.375	ng	
90) Benzo(a)pyrene	15.41	252	100697	4.515	ng	99
92) Benzo(a,h,i)perylene	17.39	276	50203	2.961	ng	96

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

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