

Data Path : Z:\HPCHEM1\BNA F\DATA\BF082417\
 Data File : BF098006.D
 Acq On : 24 Aug 2017 17:13
 Operator : SJ/JU
 Sample : I4919-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SASP2P-TP-106-W(0-5)COMP

Manual Integrations
 APPROVED

Sohil
 8/25/2017 5:40:03 PM

Quant Time: Aug 25 01:42:49 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 21 15:59:04 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	122848	20.00	ng	-0.01
21) Naphthalene-d8	8.02	136	479604	20.00	ng	-0.01
38) Acenaphthene-d10	9.77	164	216723	20.00	ng	-0.01
63) Phenanthrene-d10	11.25	188	419160	20.00	ng	-0.01
75) Chrysene-d12	13.89	240	314125	20.00	ng	-0.01
86) Perylene-d12	15.30	264	271755	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1077062	133.36	ng	0.00
7) Phenol-d6	6.38	99	1466224	133.61	ng	0.00
23) Nitrobenzene-d5	7.30	82	896477	101.54	ng	-0.02
41) 2,4,6-Tribromophenol	10.56	330	267907	142.47	ng	-0.01
44) 2-Fluorobiphenyl	9.10	172	1130032	76.61	ng	-0.01
78) Terphenyl-d14	12.84	244	1003869	66.64	ng	0.00
Target Compounds						
10) Phenol	6.39	94	33572	2.616	ng	99
49) Dimethylphthalate	9.49	163	211682	13.362	ng	99
70) Phenanthrene	11.27	178	298757	13.376	ng	99
71) Anthracene	11.33	178	49523	2.186	ng	98
74) Fluoranthene	12.46	202	288455	12.373	ng	97
77) Pyrene	12.69	202	261166	10.165	ng	98
80) Benzo(a)anthracene	13.87	228	99207	5.269	ng	99
82) Chrysene	13.91	228	101915	5.442	ng	97
85) Indeno(1,2,3-cd)pyrene	16.66	276	42764	3.104	ng	94
87) Benzo(b)fluoranthene	14.90	252	96192m	5.616	ng	
88) Benzo(k)fluoranthene	14.92	252	38705m	2.405	ng	
89) Benzo(a)pyrene	15.24	252	74997	4.946	ng	96
91) Benzo(a,h,i)perylene	17.07	276	42405	3.292	ng	95

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

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