

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110518\
 Data File : BF110479.D
 Acq On : 5 Nov 2018 21:39
 Operator : JU/SJ
 Sample : J5816-07
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-4A-S

Manual Integrations
 APPROVED

Sohil
 11/6/2018 9:01:19 AM

Quant Time: Nov 06 04:52:45 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 16:06:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	107580	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	427626	20.00	ng	0.00
39) Acenaphthene-d10	10.00	164	171745	20.00	ng	0.00
64) Phenanthrene-d10	11.49	188	253691	20.00	ng	0.00
76) Chrysene-d12	14.13	240	184084	20.00	ng	0.00
87) Perylene-d12	15.66	264	140361	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	539543	81.67	ng	0.00
7) Phenol-d6	6.58	99	689522	81.60	ng	0.00
23) Nitrobenzene-d5	7.52	82	417088	57.85	ng	0.00
42) 2,4,6-Tribromophenol	10.79	330	105655	62.10	ng	0.00
45) 2-Fluorobiphenyl	9.31	172	531677	48.61	ng	0.00
79) Terphenyl-d14	13.07	244	308324	34.83	ng	0.00
Target Compounds						
10) Phenol	6.59	94	52507	5.813	ng	Qvalue # 64
50) Dimethylphthalate	9.70	163	122099	9.631	ng	99
71) Phenanthrene	11.51	178	78093	5.883	ng	97
75) Fluoranthene	12.70	202	192581	14.295	ng	98
78) Pyrene	12.94	202	149719	11.345	ng	98
81) Benzo(a)anthracene	14.12	228	86616	7.934	ng	99
83) Chrysene	14.16	228	92855	8.955	ng	97
86) Indeno(1,2,3-cd)pyrene	17.23	276	28380	3.299	ng	96
88) Benzo(b)fluoranthene	15.20	252	94634m	11.676	ng	
89) Benzo(k)fluoranthene	15.23	252	37864m	4.752	ng	
90) Benzo(a)pyrene	15.59	252	47417	6.358	ng	# 98
92) Benzo(a,h,i)perylene	17.70	276	22307	3.518	ng	99

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

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