

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112619\
 Data File : BF118015.D
 Acq On : 26 Nov 2019 18:56
 Operator : JU
 Sample : K5948-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-1-0-6

Manual Integrations
 APPROVED

jagrut
 11/27/2019 11:33:42 PM

Quant Time: Nov 27 01:04:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 26 05:38:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	110623	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	416380	20.00	ng	0.00
39) Acenaphthene-d10	9.95	164	217009	20.00	ng	0.00
64) Phenanthrene-d10	11.44	188	287885	20.00	ng	0.00
76) Chrysene-d12	14.09	240	250777	20.00	ng	0.00
87) Perylene-d12	15.60	264	206624	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	785617	125.71	ng	0.02
7) Phenol-d6	6.53	99	855227	113.46	ng	0.00
23) Nitrobenzene-d5	7.46	82	505991	72.24	ng	0.00
42) 2,4,6-Tribromophenol	10.74	330	218215	94.98	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	1038275	85.84	ng	0.00
79) Terphenyl-d14	13.03	244	842060	61.37	ng	0.00
Target Compounds						
10) Phenol	6.53	94	17771	2.269	ng	Qvalue # 74
50) Dimethylphthalate	9.65	163	187332	12.310	ng	98
71) Phenanthrene	11.46	178	50312	3.476	ng	97
78) Pyrene	12.89	202	85522	4.220	ng	98
81) Benzo(a)anthracene	14.08	228	47883	2.889	ng	92
83) Chrysene	14.12	228	45572	2.838	ng	99
84) Bis(2-ethylhexyl)phthalate	14.06	149	21942	2.079	ng	97
88) Benzo(b)fluoranthene	15.15	252	48600m	3.620	ng	
90) Benzo(a)pyrene	15.53	252	34274	2.903	ng	# 88

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

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