

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082718\
 Data File : BF108513.D
 Acq On : 27 Aug 2018 16:28
 Operator : JU/SJ
 Sample : J4646-15
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP1-Q4-FD01

Manual Integrations
 APPROVED

Sohil
 8/27/2018 5:20:43 PM

Quant Time: Aug 27 17:05:45 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 22 11:01:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.00	152	115644	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	451054	20.00	ng	0.00
38) Acenaphthene-d10	10.04	164	215705	20.00	ng	-0.01
63) Phenanthrene-d10	11.54	188	354179	20.00	ng	-0.01
75) Chrysene-d12	14.19	240	278878	20.00	ng	-0.01
86) Perylene-d12	15.75	264	269016	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	446640	69.92	ng	0.00
7) Phenol-d6	6.62	99	610630	71.94	ng	-0.01
23) Nitrobenzene-d5	7.56	82	402622	49.81	ng	-0.01
41) 2,4,6-Tribromophenol	10.84	330	140977	65.52	ng	-0.01
44) 2-Fluorobiphenyl	9.35	172	791636	58.92	ng	-0.01
78) Terphenyl-d14	13.12	244	576191	52.53	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.74	163	118067	7.858	ng	100
70) Phenanthrene	11.56	178	47344	2.834	ng	98
74) Fluoranthene	12.76	202	75281	4.129	ng	95
77) Pyrene	12.99	202	70127	3.972	ng	98
80) Benzo(a)anthracene	14.18	228	35669	2.229	ng	100
82) Chrysene	14.22	228	34095	2.324	ng	96
87) Benzo(b)fluoranthene	15.28	252	48942m	3.045	ng	
89) Benzo(a)pyrene	15.68	252	31844	2.212	ng	94

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

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