

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082519\
 Data File : BF116530.D
 Acq On : 25 Aug 2019 14:26
 Operator : HP/JU
 Sample : K4475-07
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 T1-2-D

Manual Integrations
 APPROVED

mohammad
 8/29/2019 7:53:05 AM

Quant Time: Aug 26 07:17:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 23 13:38:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	150343	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	570903	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	273169	20.00	ng	0.00
64) Phenanthrene-d10	11.40	188	404961	20.00	ng	0.00
76) Chrysene-d12	14.03	240	213075	20.00	ng	0.00
87) Perylene-d12	15.50	264	228697	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	854492	83.05	ng	0.01
7) Phenol-d6	6.52	99	1146092	90.84	ng	0.01
23) Nitrobenzene-d5	7.44	82	734988	70.43	ng	0.00
42) 2,4,6-Tribromophenol	10.70	330	166358	70.92	ng	0.00
45) 2-Fluorobiphenyl	9.24	172	1138248	67.10	ng	0.00
79) Terphenyl-d14	12.98	244	821709	72.03	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.63	163	142444	7.474	ng	99
71) Phenanthrene	11.42	178	48812	2.213	ng	97
75) Fluoranthene	12.60	202	87464	3.835	ng	99
78) Pyrene	12.83	202	75130	4.409	ng	99
81) Benzo(a)anthracene	14.02	228	32655	2.366	ng	97
83) Chrysene	14.05	228	28655m	2.049	ng	
88) Benzo(b)fluoranthene	15.07	252	38623m	2.734	ng	
90) Benzo(a)pyrene	15.43	252	25440	2.037	ng	# 80

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

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