

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102419\
 Data File : BF117524.D
 Acq On : 24 Oct 2019 21:32
 Operator : JU
 Sample : K5499-13 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 707-58-(0-1)

Manual Integrations
 APPROVED

mohammad
 10/26/2019 10:03:17 AM

Quant Time: Oct 25 04:33:59 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 18:53:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	50091	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	204353	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	93937	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	134225	20.00	ng	0.00
76) Chrysene-d12	13.90	240	98633	20.00	ng	0.00
87) Perylene-d12	15.33	264	78950	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	118345	35.15	ng	0.00
7) Phenol-d6	6.39	99	168711	37.31	ng	0.00
23) Nitrobenzene-d5	7.30	82	96511	23.32	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	24770	30.48	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	173116	25.97	ng	0.00
79) Terphenyl-d14	12.83	244	130445	21.57	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.49	163	20251	3.024	ng	Qvalue 100
71) Phenanthrene	11.27	178	19632	2.519	ng	99
75) Fluoranthene	12.47	202	38357	4.916	ng	98
78) Pyrene	12.70	202	32873	3.880	ng	99
81) Benzo(a)anthracene	13.89	228	16868	2.535	ng	98
83) Chrysene	13.92	228	15092	2.315	ng	94
88) Benzo(b)fluoranthene	14.92	252	17851m	3.829	ng	
90) Benzo(a)pyrene	15.27	252	10845	2.559	ng	# 86

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

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