

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062719\
 Data File : BF115238.D
 Acq On : 27 Jun 2019 16:07
 Operator : HP/JU
 Sample : K3528-06
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-08-COMP

Manual Integrations
 APPROVED

mohammad
 6/28/2019 2:58:24 PM

Quant Time: Jun 28 05:45:01 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF062119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 28 05:26:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.60	152	228937	20.00	ng	0.00
21) Naphthalene-d8	7.88	136	859574	20.00	ng	0.00
39) Acenaphthene-d10	9.62	164	414058	20.00	ng	0.00
64) Phenanthrene-d10	11.10	188	811712	20.00	ng	0.00
76) Chrysene-d12	13.71	240	661766	20.00	ng	0.00
87) Perylene-d12	15.07	264	553365	20.00	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.20	112	1323834	89.23	ng	0.00
7) Phenol-d6	6.24	99	1667511	92.61	ng	0.00
23) Nitrobenzene-d5	7.16	82	984552	70.46	ng	0.00
42) 2,4,6-Tribromophenol	10.41	330	431756	98.88	ng	0.00
45) 2-Fluorobiphenyl	8.95	172	1748081	71.71	ng	0.00
79) Terphenyl-d14	12.68	244	1964521	63.12	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
50) Dimethylphthalate	9.35	163	238590	7.832	ng	99
71) Phenanthrene	11.12	178	287496	6.578	ng	96
75) Fluoranthene	12.30	202	462396	10.604	ng	96
78) Pyrene	12.52	202	394667	7.760	ng	99
81) Benzo(a)anthracene	13.70	228	209800	4.418	ng	98
83) Chrysene	13.74	228	170897	3.929	ng	96
88) Benzo(b)fluoranthene	14.70	252	226954m	6.573	ng	
90) Benzo(a)pyrene	15.01	252	130403	4.190	ng	99
92) Benzo(g,h,i)perylene	16.70	276	61747	2.100	ng	97

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

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