

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP050520\
 Data File : BP002246.D
 Acq On : 05 May 2020 20:27
 Operator : CG/JU
 Sample : L2483-02 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 6FT-TP-1

Manual Integrations
 APPROVED

mohammad
 5/6/2020 6:44:39 PM

Quant Time: May 06 01:45:57 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP050420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 05 13:34:25 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	263974	20.00	ng	0.00
21) Naphthalene-d8	10.39	136	1211015	20.00	ng	0.00
39) Acenaphthene-d10	14.26	164	762596	20.00	ng	0.00
64) Phenanthrene-d10	17.01	188	1663439	20.00	ng	0.00
76) Chrysene-d12	21.11	240	1768095	20.00	ng	0.00
87) Perylene-d12	23.31	264	1798649	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	344003	21.50	ng	0.00
7) Phenol-d6	6.79	99	510937	21.74	ng	0.00
23) Nitrobenzene-d5	8.75	82	302563	12.90	ng	0.00
42) 2,4,6-Tribromophenol	15.75	330	172061	21.32	ng	0.00
45) 2-Fluorobiphenyl	12.87	172	815019	14.97	ng	0.00
79) Terphenyl-d14	19.59	244	1294572	17.94	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	13.72	163	125159	2.082	ng	98
71) Phenanthrene	17.05	178	313621	3.353	ng	100
75) Fluoranthene	19.03	202	537796	5.011	ng	98
78) Pyrene	19.38	202	487259	4.057	ng	99
81) Benzo(a)anthracene	21.09	228	282023	2.459	ng	98
83) Chrysene	21.14	228	242629m	2.154	ng	
88) Benzo(b)fluoranthene	22.65	252	324236m	2.998	ng	
90) Benzo(a)pyrene	23.22	252	233434	2.242	ng	97

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

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