

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP052920\
 Data File : BP002327.D
 Acq On : 29 May 2020 22:01
 Operator : CG/JU
 Sample : L2744-11 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 NM79-COMP-2020-0-6

Quant Time: May 30 02:30:06 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP052720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 29 15:09:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	380794	20.00	ng	0.00
21) Naphthalene-d8	10.38	136	1579266	20.00	ng	0.00
39) Acenaphthene-d10	14.25	164	929964	20.00	ng	0.00
64) Phenanthrene-d10	17.01	188	1916463	20.00	ng	0.00
76) Chrysene-d12	21.11	240	1707031	20.00	ng	0.00
86) Perylene-d12	23.32	264	1770368	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	1117463	55.11	ng	0.00
7) Phenol-d6	6.79	99	1533961	55.62	ng	0.00
23) Nitrobenzene-d5	8.75	82	984133	37.47	ng	0.00
42) 2,4,6-Tribromophenol	15.75	330	498104	56.51	ng	0.00
45) 2-Fluorobiphenyl	12.87	172	2147206	40.74	ng	0.00
79) Terphenyl-d14	19.59	244	2890731	41.01	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.71	163	328737	5.222	ng	99
75) Fluoranthene	19.03	202	433318	4.144	ng	99
78) Pyrene	19.39	202	419074	3.943	ng	99
81) Benzo(a)anthracene	21.10	228	233708	2.392	ng	97
83) Chrysene	21.14	228	222238	2.370	ng	97
88) Benzo(b)fluoranthene	22.66	252	308014	3.191	ng	# 96
90) Benzo(a)pyrene	23.22	252	223840	2.500	ng	# 94

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

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