

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081920\
 Data File : BP003043.D
 Acq On : 19 Aug 2020 21:12
 Operator : CG/JU
 Sample : L3712-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CHRT-26572

Quant Time: Aug 20 02:03:24 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 17 14:07:35 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	344305	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	1417407	20.00	ng	-0.01
39) Acenaphthene-d10	14.32	164	849751	20.00	ng	-0.01
64) Phenanthrene-d10	17.07	188	1657584	20.00	ng	-0.01
76) Chrysene-d12	21.16	240	1478018	20.00	ng	-0.01
86) Perylene-d12	23.41	264	1625085	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	1675406	74.47	ng	0.00
7) Phenol-d6	6.86	99	2034468	65.14	ng	0.00
23) Nitrobenzene-d5	8.82	82	1173285	44.83	ng	-0.01
42) 2,4,6-Tribromophenol	15.81	330	896620	84.90	ng	-0.01
45) 2-Fluorobiphenyl	12.94	172	3304747	53.14	ng	-0.01
79) Terphenyl-d14	19.65	244	3915447	53.03	ng	0.00
Target Compounds						
10) Phenol	6.89	94	68280	2.171	ng	91
50) Dimethylphthalate	13.77	163	420894	6.333	ng	99
75) Fluoranthene	19.09	202	249833	2.231	ng	97
78) Pyrene	19.44	202	335980	3.252	ng	97
84) Bis(2-ethylhexyl)phthalate	21.10	149	235756	4.123	ng	99
88) Benzo(b)fluoranthene	22.73	252	292264	2.549	ng	98

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

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