

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP072320\
 Data File : BP002849.D
 Acq On : 23 Jul 2020 20:10
 Operator : CG/JU
 Sample : L3402-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CHASSIE-WASH

Quant Time: Jul 24 01:23:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 13 10:54:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	48151	20.00	ng	-0.01
21) Naphthalene-d8	10.31	136	203440	20.00	ng	-0.02
39) Acenaphthene-d10	14.19	164	132916	20.00	ng	-0.02
64) Phenanthrene-d10	16.96	188	313525	20.00	ng	-0.01
76) Chrysene-d12	21.07	240	421514	20.00	ng	0.00
86) Perylene-d12	23.25	264	521011	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.17	112	365582	128.10	ng	-0.01
7) Phenol-d6	6.75	99	444673	109.18	ng	-0.01
23) Nitrobenzene-d5	8.69	82	352225	92.60	ng	-0.02
42) 2,4,6-Tribromophenol	15.70	330	210613	151.80	ng	-0.01
45) 2-Fluorobiphenyl	12.80	172	754427	87.81	ng	-0.02
79) Terphenyl-d14	19.55	244	1324466	72.04	ng	-0.01
Target Compounds						
6) Aniline	6.89	93	13049	2.531	ng	99
10) Phenol	6.78	94	13254	3.177	ng	# 62
20) 3+4-Methylphenols	8.33	107	92816	23.615	ng	94
31) Naphthalene	10.36	128	23361	2.175	ng	98
32) Benzoic acid	9.69	122	13399	12.176	ng	85
50) Dimethylphthalate	13.66	163	159020	15.688	ng	100

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

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