

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100720\
 Data File : BP003565.D
 Acq On : 07 Oct 2020 22:15
 Operator : CG/JU
 Sample : L4260-12
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PAV-B23-100220-0-10

Quant Time: Oct 08 01:54:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 05 14:33:14 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.493	152	83416	20.00	ng	0.00
21) Naphthalene-d8	10.263	136	327915	20.00	ng	#-0.01
39) Acenaphthene-d10	14.151	164	188373	20.00	ng	0.00
64) Phenanthrene-d10	16.910	188	371780	20.00	ng	0.00
76) Chrysene-d12	21.016	240	327831	20.00	ng	-0.01
86) Perylene-d12	23.186	264	382411	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.111	112	538592	112.74	ng	0.00
7) Phenol-d6	6.699	99	605320	95.06	ng	0.00
23) Nitrobenzene-d5	8.646	82	596645	106.89	ng	0.00
42) 2,4,6-Tribromophenol	15.657	330	322940	112.40	ng	0.00
45) 2-Fluorobiphenyl	12.763	172	1200718	93.22	ng	-0.01
79) Terphenyl-d14	19.492	244	1488177	94.58	ng	-0.01
Target Compounds						
						Qvalue
10) Phenol	6.728	94	28021	4.61	ng	# 58
17) 2-Methylphenol	7.987	107	12330	2.80	ng	# 89
20) 3+4-Methylphenols	8.334	107	33007	5.64	ng	90
27) 2,4-Dimethylphenol	9.516	122	13456	3.12	ng	90
31) Naphthalene	10.316	128	178820	10.32	ng	99

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

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