

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
 Data File : BP013057.D
 Acq On : 11 Dec 2022 00:58
 Operator : CG/JU
 Sample : N5927-21
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXZA7

Quant Time: Dec 12 03:30:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 12 01:32:54 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.969	152	615498	20.000 ng/u1	0.00
20) Naphthalene-d8	10.781	136	2691340	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.610	164	1830335	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.363	188	4164213	20.000 ng/u1	0.00
79) Chrysene-d12	21.451	240	3850569	20.000 ng/u1	0.00
88) Perylene-d12	23.939	264	3850669	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.352	96	70982	4.931 ng/uL	0.00
4) Pyridine-d5	3.781	84	357400	8.740 ng/u1	0.00
7) Phenol-d5	7.122	99	392688	7.833 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	1177690	38.941 ng/u1	0.00
11) 2-Chlorophenol-d4	7.499	132	1178957	29.873 ng/u1	0.00
15) 4-Methylphenol-d8	8.675	113	780604	19.003 ng/u1	0.00
21) Nitrobenzene-d5	9.140	128	779520	39.421 ng/u1	0.00
24) 2-Nitrophenol-d4	9.863	143	880471	39.551 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	1469409	33.982 ng/u1	0.00
31) 4-Chloroaniline-d4	10.922	131	1827748	29.942 ng/u1	0.00
46) Dimethylphthalate-d6	14.016	166	5618369	39.940 ng/u1	0.00
49) Acenaphthylene-d8	14.304	160	6080839	39.340 ng/u1	0.00
54) 4-Nitrophenol-d4	14.798	143	152173	6.014 ng/u1	0.00
60) Fluorene-d10	15.604	176	4813856	40.450 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	943901	35.223 ng/u1	0.00
73) Anthracene-d10	17.463	188	7815021	41.599 ng/u1	0.00
81) Pyrene-d10	19.698	212	9170167	41.490 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.780	264	8121516	42.314 ng/u1	0.00

Target Compounds Qvalue

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

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