

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012476.D
 Acq On : 03 Nov 2022 06:19
 Operator : CG/JU
 Sample : PB148676BL
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK676

Quant Time: Nov 03 07:25:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 03 00:09:04 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.781	152	324972	20.000 ng/u1	0.00
20) Naphthalene-d8	10.581	136	1490160	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.434	164	1033769	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.198	188	2355504	20.000 ng/u1	0.00
79) Chrysene-d12	21.292	240	2213351	20.000 ng/u1	0.00
88) Perylene-d12	23.674	264	1973199	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.234	96	41698	5.167 ng/uL	0.00
4) Pyridine-d5	3.652	84	350750	15.129 ng/u1	0.00
7) Phenol-d5	6.958	99	808563	27.740 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	468168	26.906 ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	618559	29.060 ng/u1	0.00
15) 4-Methylphenol-d8	8.499	113	669426	29.039 ng/u1	0.00
21) Nitrobenzene-d5	8.952	128	326263	33.763 ng/u1	0.00
24) 2-Nitrophenol-d4	9.669	143	374848	37.373 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	665538	30.618 ng/u1	0.00
31) 4-Chloroaniline-d4	10.734	131	587397	18.149 ng/u1	0.00
46) Dimethylphthalate-d6	13.851	166	2156182	30.469 ng/u1	0.00
49) Acenaphthylene-d8	14.128	160	2383018	29.467 ng/u1	0.00
54) 4-Nitrophenol-d4	14.663	143	372945	27.845 ng/u1	0.00
60) Fluorene-d10	15.434	176	1856244	30.637 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	235953	19.240 ng/u1	0.00
73) Anthracene-d10	17.292	188	3072054	30.269 ng/u1	-0.01
81) Pyrene-d10	19.539	212	3535738	31.818 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.521	264	2966721	30.771 ng/u1	0.00

Target Compounds Qvalue

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

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