

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
 Data File : BP012829.D
 Acq On : 28 Nov 2022 22:47
 Operator : CG/JU
 Sample : PB149199BL
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK199

Quant Time: Nov 29 00:46:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Nov 27 22:55:08 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	8.016	152	743051	20.000 ng/u1	0.00
20) Naphthalene-d8	10.828	136	3213777	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.651	164	1963748	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.404	188	3347919	20.000 ng/u1	0.00
79) Chrysene-d12	21.486	240	2163568	20.000 ng/u1	-0.01
88) Perylene-d12	23.998	264	2508995	20.000 ng/u1	-0.01

System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.399	96	95769	5.523 ng/uL	0.00
4) Pyridine-d5	3.822	84	1264061	24.359 ng/u1	0.00
7) Phenol-d5	7.163	99	1576361	25.817 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	1022336	27.172 ng/u1	0.00
11) 2-Chlorophenol-d4	7.540	132	1287805	27.680 ng/u1	0.00
15) 4-Methylphenol-d8	8.716	113	1258832	25.588 ng/u1	0.00
21) Nitrobenzene-d5	9.181	128	608617	30.941 ng/u1	0.00
24) 2-Nitrophenol-d4	9.904	143	672133	32.181 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	1266186	25.982 ng/u1	0.00
31) 4-Chloroaniline-d4	10.963	131	1761973	26.403 ng/u1	0.00
46) Dimethylphthalate-d6	14.057	166	3682920	27.247 ng/u1	0.00
49) Acenaphthylene-d8	14.345	160	4309824	27.327 ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	432530	17.948 ng/u1	0.00
60) Fluorene-d10	15.639	176	3182253	26.883 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	346949	21.682 ng/u1	0.00
73) Anthracene-d10	17.504	188	4240888	28.777 ng/u1	0.00
81) Pyrene-d10	19.727	212	3833826	30.496 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.833	264	3572826	28.991 ng/u1	-0.01

Target Compounds	Qvalue

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

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