

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
 Data File : BP022185.D
 Acq On : 03 Oct 2024 07:42
 Operator : RC/JU
 Sample : PB163456BL
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK456

Quant Time: Oct 03 08:57:02 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 24 15:23:13 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.699	152	116854	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.457	136	460727	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.304	164	348969	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.104	188	842069	20.000	ng/u1	-0.01
79) Chrysene-d12	21.539	240	913780	20.000	ng/u1	0.00
88) Perylene-d12	24.821	264	1035975	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	14045	4.708	ng/uL	0.00
4) Pyridine-d5	3.634	84	198776	23.310	ng/u1	0.00
7) Phenol-d5	6.881	99	253403	26.510	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.040	67	148817	25.284	ng/u1	0.00
11) 2-Chlorophenol-d4	7.240	132	200433	28.981	ng/u1	0.00
15) 4-Methylphenol-d8	8.399	113	212066	27.998	ng/u1	0.00
21) Nitrobenzene-d5	8.834	128	99610	28.886	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.552	143	120773	30.889	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.081	165	236707	27.877	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.593	131	280479	27.685	ng/u1	-0.01
46) Dimethylphthalate-d6	13.710	166	798247	29.591	ng/u1	-0.01
49) Acenaphthylene-d8	13.992	160	840120	29.415	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.522	143	108996	23.622	ng/u1	0.00
60) Fluorene-d10	15.310	176	677183	28.380	ng/u1	-0.02
65) 4,6-Dinitro-2-methylph...	15.445	200	105145	20.112	ng/u1	0.00
73) Anthracene-d10	17.210	188	1153385	29.340	ng/u1	0.00
81) Pyrene-d10	19.586	212	1450954	30.649	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.586	264	1589950	29.179	ng/u1	-0.02

Target Compounds Qvalue

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

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