

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
 Data File : BP022158.D
 Acq On : 02 Oct 2024 12:04
 Operator : RC/JU
 Sample : PB163456BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK456

Quant Time: Oct 02 13:40:28 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	121059	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.457	136	468923	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.304	164	364636	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.104	188	886146	20.000	ng/u1	-0.01
79) Chrysene-d12	21.539	240	1009781	20.000	ng/u1	0.00
88) Perylene-d12	24.816	264	1136214	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	15332	4.961	ng/uL	-0.01
4) Pyridine-d5	3.628	84	214015	24.225	ng/u1	0.00
7) Phenol-d5	6.881	99	260855	26.342	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.034	67	160518	26.324	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.234	132	200881	28.037	ng/u1	-0.01
15) 4-Methylphenol-d8	8.399	113	210728	26.855	ng/u1	0.00
21) Nitrobenzene-d5	8.834	128	102106	29.092	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.552	143	125565	31.553	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.087	165	231784	26.820	ng/u1	0.00
31) 4-Chloroaniline-d4	10.587	131	271672	26.347	ng/u1	-0.02
46) Dimethylphthalate-d6	13.710	166	834810	29.617	ng/u1	-0.01
49) Acenaphthylene-d8	13.993	160	841670	28.203	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.522	143	119317	24.748	ng/u1	0.00
60) Fluorene-d10	15.310	176	692291	27.766	ng/u1	-0.02
65) 4,6-Dinitro-2-methylph...	15.434	200	148595	27.009	ng/u1	-0.01
73) Anthracene-d10	17.198	188	1194198	28.867	ng/u1	-0.02
81) Pyrene-d10	19.581	212	1507259	28.811	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.586	264	1685982	28.212	ng/u1	-0.02

Target Compounds Qvalue

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

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