

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100424\
 Data File : BP022244.D
 Acq On : 05 Oct 2024 01:58
 Operator : RC/JU
 Sample : P4264-12
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCY01

Quant Time: Oct 05 03:30:33 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 04 05:25:28 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.687	152	115026	20.000	ng/u1	0.00
20) Naphthalene-d8	10.446	136	437009	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.304	164	348821	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.104	188	831617	20.000	ng/u1	0.01
79) Chrysene-d12	21.527	240	910018	20.000	ng/u1	0.00
88) Perylene-d12	24.804	264	1074310	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	11708	3.987	ng/uL	0.00
4) Pyridine-d5	3.634	84	38783	4.620	ng/u1	0.00
7) Phenol-d5	6.875	99	67229	7.145	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.028	67	139982	24.160	ng/u1	0.00
11) 2-Chlorophenol-d4	7.234	132	161061	23.659	ng/u1	0.00
15) 4-Methylphenol-d8	8.387	113	123878	16.615	ng/u1	-0.01
21) Nitrobenzene-d5	8.828	128	88441	27.039	ng/u1	0.00
24) 2-Nitrophenol-d4	9.540	143	105479	28.441	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	210975	26.195	ng/u1	0.00
31) 4-Chloroaniline-d4	10.581	131	206833	21.523	ng/u1	0.00
46) Dimethylphthalate-d6	13.704	166	785337	29.125	ng/u1	0.00
49) Acenaphthylene-d8	13.993	160	761019	26.657	ng/u1	0.00
54) 4-Nitrophenol-d4	14.522	143	25936	5.623	ng/u1	0.00
60) Fluorene-d10	15.316	176	683411	28.653	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.434	200	123151	23.852	ng/u1	0.00
73) Anthracene-d10	17.198	188	1266831	32.631	ng/u1	0.00
81) Pyrene-d10	19.581	212	1608522	34.118	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.586	264	1849591	32.733	ng/u1	0.01

Target Compounds Qvalue

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

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