

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
 Data File : BP022764.D
 Acq On : 31 Oct 2024 10:44
 Operator : RC/JU
 Sample : P4592-10
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 F7H18

Quant Time: Oct 31 11:40:46 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 30 10:56:39 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.634	152	81986	20.000	ng/u1	0.00
20) Naphthalene-d8	10.393	136	315991	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.263	164	243995	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.063	188	605764	20.000	ng/u1	0.00
79) Chrysene-d12	21.486	240	719205	20.000	ng/u1	#-0.01
88) Perylene-d12	24.733	264	895469	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	5740	2.721	ng/uL	0.00
4) Pyridine-d5	3.593	84	77755	13.425	ng/u1	0.00
7) Phenol-d5	6.834	99	114290	16.561	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.981	67	63738	15.714	ng/u1	0.00
11) 2-Chlorophenol-d4	7.181	132	82020	16.749	ng/u1	0.00
15) 4-Methylphenol-d8	8.340	113	95783	17.391	ng/u1	0.00
21) Nitrobenzene-d5	8.775	128	39056	16.075	ng/u1	0.00
24) 2-Nitrophenol-d4	9.493	143	49814	17.709	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.034	165	105179	17.593	ng/u1	0.00
31) 4-Chloroaniline-d4	10.534	131	116837	16.539	ng/u1	0.00
46) Dimethylphthalate-d6	13.657	166	359962	18.568	ng/u1	0.00
49) Acenaphthylene-d8	13.945	160	378916	18.403	ng/u1	0.00
54) 4-Nitrophenol-d4	14.522	143	37747	11.606	ng/u1	0.01
60) Fluorene-d10	15.281	176	323764	19.255	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.416	200	28867	7.246	ng/u1	0.01
73) Anthracene-d10	17.163	188	568589	19.953	ng/u1	0.00
81) Pyrene-d10	19.545	212	735908	19.349	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.515	264	970064	20.285	ng/u1	0.00

Target Compounds	Qvalue

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

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