

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012423.D
 Acq On : 01 Nov 2022 12:41
 Operator : CG/JU
 Sample : N5270-07
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHA89

Quant Time: Nov 01 22:57:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 01 01:03:58 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	321819	20.000	ng/u1	0.00
20) Naphthalene-d8	10.593	136	1456138	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.445	164	936451	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.204	188	2036933	20.000	ng/u1	0.00
79) Chrysene-d12	21.298	240	2194606	20.000	ng/u1	0.00
88) Perylene-d12	23.686	264	2141669	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	38621	4.832	ng/uL	0.00
4) Pyridine-d5	3.664	84	168200	7.326	ng/u1	0.00
7) Phenol-d5	6.963	99	224420	7.775	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	609764	35.387	ng/u1	0.00
11) 2-Chlorophenol-d4	7.322	132	587443	27.869	ng/u1	0.00
15) 4-Methylphenol-d8	8.510	113	438884	19.225	ng/u1	0.00
21) Nitrobenzene-d5	8.957	128	414387	43.884	ng/u1	0.00
24) 2-Nitrophenol-d4	9.681	143	408929	41.723	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	733282	34.523	ng/u1	0.00
31) 4-Chloroaniline-d4	10.740	131	857805	27.123	ng/u1	0.00
46) Dimethylphthalate-d6	13.857	166	2747344	42.857	ng/u1	0.00
49) Acenaphthylene-d8	14.139	160	3065316	41.843	ng/u1	0.00
54) 4-Nitrophenol-d4	14.669	143	80936	6.671	ng/u1	0.00
60) Fluorene-d10	15.439	176	2368764	43.158	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	393260	37.082	ng/u1	0.00
73) Anthracene-d10	17.304	188	3872342	44.121	ng/u1	0.00
81) Pyrene-d10	19.545	212	4633825	42.055	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.533	264	4658894	44.521	ng/u1	-0.01

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

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