

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102721\
 Data File : BG050772.D
 Acq On : 28 Oct 2021 11:37
 Operator : CG/JU
 Sample : M4215-19
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JDGF6

Quant Time: Oct 28 12:13:58 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102721.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 26 18:31:25 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.250	152	35431	20.000	ng/u1	0.00
20) Naphthalene-d8	11.076	136	158739	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.872	164	102679	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.621	188	213046	20.000	ng/u1	0.00
79) Chrysene-d12	21.928	240	196267	20.000	ng/u1	# 0.00
88) Perylene-d12	25.348	264	199440	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.602	96	4138	4.602	ng/uL	0.00
4) Pyridine-d5	4.031	84	22145	7.777	ng/u1	0.00
7) Phenol-d5	7.386	99	114111	34.658	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.557	67	73152	34.533	ng/u1	0.00
11) 2-Chlorophenol-d4	7.774	132	82113	35.505	ng/u1	0.00
15) 4-Methylphenol-d8	8.943	113	91561	36.018	ng/u1	0.00
21) Nitrobenzene-d5	9.419	128	43577	35.965	ng/u1	0.00
24) 2-Nitrophenol-d4	10.148	143	50892	37.247	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.688	165	81864	34.116	ng/u1	0.00
31) 4-Chloroaniline-d4	11.205	131	101940	28.947	ng/u1	0.00
46) Dimethylphthalate-d6	14.266	166	265383	36.597	ng/u1	0.00
49) Acenaphthylene-d8	14.572	160	339179	37.663	ng/u1	0.00
54) 4-Nitrophenol-d4	15.054	143	45170	34.418	ng/u1	0.00
60) Fluorene-d10	15.865	176	239513	38.434	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.976	200	36785	29.505	ng/u1	0.00
73) Anthracene-d10	17.721	188	358409	39.237	ng/u1	0.00
81) Pyrene-d10	19.995	212	399347	35.705	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.118	264	358286	34.916	ng/u1	0.00

Target Compounds Qvalue

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

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